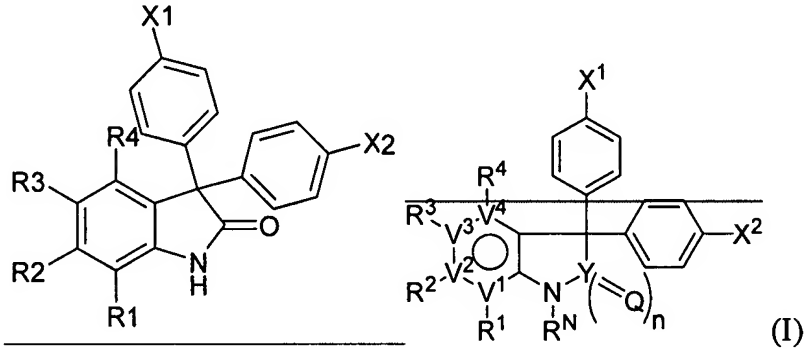
AMENDMENTS TO THE CLAIMS

The following listing of claims replaces all prior versions, and listings, of claims in this application.

Claim 1 (**currently amended**): A method of treating a mammal suffering from or being susceptible to cancer, the method comprising administering to the mammal a therapeutically effective amount of a compound of the general formula (I)



(I)

wherein

~~V^1, V^2, V^3, and V^4 independently are selected from a carbon atom, a non-quaternary nitrogen atom, an oxygen atom, and a sulfur atom, and where V^4 further may be selected from a bond, so that V^1-V^2-V^3-V^4 together with the atoms to which V^1 and V^4 are attached form an aromatic or heteroaromatic ring;~~

R^1, R^2, R^3, and R^4~~, when attached to a carbon atom,~~ independently are selected from hydrogen, optionally substituted C_{1-6}-alkyl, optionally substituted C_{2-6}-alkenyl, hydroxy, optionally substituted C_{1-6}-alkoxy, optionally substituted C_{2-6}-alkenyloxy, carboxy, optionally substituted C_{1-6}-alkoxycarbonyl, optionally substituted C_{1-6}-alkylcarbonyl, optionally substituted C_{1-6}-alkylcarbonyloxy, formyl, amino, mono- and di(C_{1-6}-alkyl)amino, carbamoyl, mono- and di(C_{1-6}-alkyl)aminocarbonyl, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylsulphonylamino, cyano, carbamido, mono- and di(C_{1-6}-alkyl)aminocarbonylamino, C_{1-6}-alkanoyloxy, C_{1-6}-alkylsulphonyl, C_{1-6}-alkylsulphinyl, aminosulfonyl, mono- and di(C_{1-6}-alkyl)aminosulfonyl, nitro, optionally substituted

C_{1-6}-alkylthio, aryl, aryloxy, arylcarbonyl, arylamino, heterocyclyl, heterocyclyloxy, heterocyclylamino, heterocyclylcarbonyl, heteroaryl, heteroaryloxy, heteroarylamino, heteroarylcarbonyl, and halogen, where any C_{1-6}-alkyl as an amino substituent is optionally substituted with hydroxy, C_{1-6}-alkoxy, amino, mono- and di(C_{1-6}-alkyl)amino, carboxy, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylaminocarbonyl, or halogen(s), and wherein any aryl, heterocyclyl and heteroaryl may be optionally substituted;

~~R^1, R^2, R^3, and R^4, when attached to a nitrogen atom, independently are selected from hydrogen, optionally substituted C_{1-6}-alkyl, hydroxy, optionally substituted C_{1-6}-alkoxy, optionally substituted C_{1-6}-alkoxycarbonyl, optionally substituted C_{1-6}-alkylcarbonyl, formyl, mono- and di(C_{1-6}-alkyl)-aminocarbonyl, amino, C_{1-6}-alkylcarbonylamino, mono- and di(C_{1-6}-alkyl)amino, C_{1-6}-alkylsulphonyl, C_{1-6}-alkylsulphinyl, aryl, aryloxy, arylcarbonyl, arylamino, heterocyclyl, heterocyclyloxy, heterocyclylcarbonyl, heterocyclylamino, heteroaryl, heteroaryloxy, heteroarylcarbonyl, and heteroarylamino; where any C_{1-6}-alkyl as an amino substituent is optionally substituted with hydroxy, C_{1-6}-alkoxy, amino, mono- and di(C_{1-6}-alkyl)amino, carboxy, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylaminocarbonyl, or halogen(s), and wherein any aryl, heterocyclyl and heteroaryl may be optionally substituted;~~

or R^1 and R^2 together with the carbon atoms to which they are attached form a ring;

with the proviso that R^1, R^2, R^3 and R^4 are not all hydrogen;

X^1 and X^2 are independently selected from halogen, hydroxy, optionally substituted C_{1-6}-alkoxy, optionally substituted C_{1-6}-alkylcarbonyloxy, amino, mono- and di(C_{1-6}-alkyl)amino, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylsulphonylamino, mono- and di(C_{1-6}-alkyl)aminocarbonylamino, C_{1-6}-alkanoyloxy, mercapto, optionally substituted C_{1-6}-alkylthio, C_{1-6}-alkylsulfonyl, mono- and di(C_{1-6}-alkyl)aminosulfonyl, aryloxy, arylamino, heterocyclyloxy, heterocyclylamino, heteroaryloxy and heteroarylamino, where any C_{1-6}-alkyl as an amino or sulphur substituent is optionally substituted with hydroxy, C_{1-6}-alkoxy, amino, mono- and di(C_{1-6}-alkyl)amino, carboxy, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylaminocarbonyl, or halogen(s), and wherein any aryl, heterocyclyl and heteroaryl may be optionally substituted;

~~$>Y(=O)_n$ is selected from >C=O, >C=S, >S=O and $>S(=O)_2$; and~~

~~R^N is selected from the group consisting of hydrogen, optionally substituted C_{1-6}-alkyl, hydroxy, optionally substituted C_{1-6}-alkoxy, optionally substituted C_{1-6}-alkoxycarbonyl, optionally substituted C_{1-6}-alkylcarbonyl, formyl, mono- and di(C_{1-6}-alkyl)aminocarbonyl, amino, C_{1-6}-alkylcarbonylamino, mono- and di(C_{1-6}-alkyl)amino, C_{1-6}-alkylsulphonyl, and C_{1-6}-alkylsulphinyl; where any C_{1-6}-alkyl as an amino substituent is optionally substituted with hydroxy, C_{1-6}-alkoxy, amino, mono- and di(C_{1-6}-alkyl)amino, carboxy, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylaminocarbonyl, or halogen(s);~~ and
pharmaceutically acceptable salts and prodrugs thereof.

Claims 2-3 (**canceled**).

Claim 4 (**currently amended**): The method according to claim 1, wherein R^1 is selected from hydrogen, halogen, C_{1-6}-alkyl, trifluoromethyl and C_{1-6}-alkoxy~~, when V^1 is a carbon atom~~.

Claim 5 (**currently amended**): The method according to claim 1, wherein R^2 is selected from hydrogen, halogen, optionally substituted aryl, optionally substituted aryloxy, and optionally substituted heteroaryl~~, when V^2 is a carbon atom~~.

Claim 6 (**currently amended**): The method according to claim 1, wherein R^3 is selected from hydrogen, optionally substituted C_{1-6}-alkoxy, halogen, cyano, optionally substituted aryl, optionally substituted aryloxy, optionally substituted heteroaryl, amino, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylsulphonylamino, and mono- and di(C_{1-6}-alkyl)aminosulfonyl~~, when V^3 is a carbon atom~~.

Claim 7 (**currently amended**): The method according to claim 1, wherein R^4 is hydrogen~~, when V^4 is a carbon atom~~.

Claim 8 (**previously presented**): The method according to claim 1, wherein X^1 and X^2 independently are selected from hydroxy, OAc, NH_2, NMe_2, NHAc, $NHSO_2Me$ and $NHCONMe_2$.

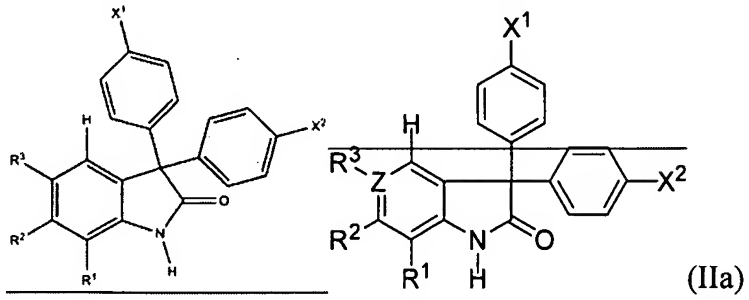
Claims 9-20 (**canceled**).

Claim 21 (**currently amended**): The method according to claim 1~~4~~, wherein R^1 is selected from fluoro, chloro, bromo, C_{1-4}-alkyl, trifluoromethyl, C_{1-4}-alkoxy, and dimethylaminocarbonyl.

Claim 22 (**canceled**).

Claim 23 (**currently amended**): The method according to claim 1~~4~~, wherein R^1 is selected from halogen, C_{1-4}-alkyl, trifluoromethyl, C_{1-4}-alkoxy, and dimethylaminocarbonyl, R^2 is selected from hydrogen and halogen, and R^3 is selected from hydrogen, halogen, C_{1-4}-alkyl, and amino; where R^2 and R^3 are not both hydrogen.

Claim 24 (**currently amended**): A method of treating a mammal suffering from or being susceptible to cancer, the method comprising administering to the mammal a therapeutically effective amount of a 3,3-diphenyl-1,3-dihydro-indol-2-one type compound of the formula (IIa)



(IIa)

wherein

R^1 is selected from hydrogen, halogen, C_{1-6}-alkyl, trifluoromethyl and C_{1-6}-alkoxy;

R^2 is selected from hydrogen, halogen, optionally substituted aryl, optionally substituted aryloxy, and optionally substituted heteroaryl;

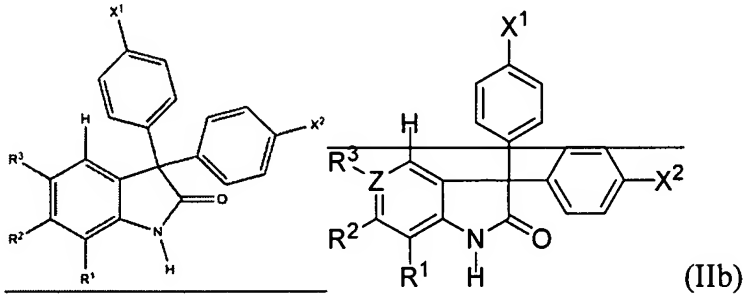
R^3 is selected from hydrogen, optionally substituted C_{1-6}-alkoxy, halogen, cyano, and optionally substituted aryl, optionally substituted aryloxy, optionally substituted heteroaryl, amino, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylsulphonylamino, and mono- and di(C_{1-6}-alkyl)aminosulfonyl;

~~Z is CH or N;~~ and

X^1 and X^2 are independently selected from halogen, OR^6, $OCOR^5$, $N(R^6)_2$, $NHCOR^5$, $NHSO_2R^5$, and $NHCON(R^6)_2$, wherein R^5 is selected from C_{1-6}-alkyl, optionally substituted aryl and optionally substituted heteroaryl, and each R^6 independently is selected from hydrogen, C_{1-6}-alkyl, optionally substituted aryl and optionally substituted heteroaryl; and

pharmaceutically acceptable salts and prodrugs thereof.

Claim 25 (**currently amended**): A method of treating a mammal suffering from or being susceptible to cancer, the method comprising administering to the mammal a therapeutically effective amount of a 3,3-diphenyl-1,3-dihydro-indol-2-one type compound of the formula (IIb)



(IIb)

wherein

R^1, R^2, and R^3~~, when attached to a carbon atom,~~ independently are selected from hydrogen, optionally substituted C_{1-6}-alkyl, optionally substituted C_{2-6}-alkenyl, hydroxy, optionally substituted C_{1-6}-alkoxy, optionally substituted C_{2-6}-alkenyloxy, carboxy, optionally substituted C_{1-6}-alkoxycarbonyl, optionally substituted C_{1-6}-alkylcarbonyl, optionally substituted C_{1-6}-alkylcarbonyloxy, formyl, amino, mono- and di(C_{1-6}-alkyl)amino, carbamoyl, mono- and di(C_{1-6}-alkyl)aminocarbonyl, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylsulphonylamino, cyano, carbamido, mono- and di(C_{1-6}-alkyl)aminocarbonylamino, C_{1-6}-alkanoyloxy, C_{1-6}-alkylsulphonyl, C_{1-6}-alkylsulphinyl, aminosulfonyl, mono- and di(C_{1-6}-alkyl)aminosulfonyl, nitro, optionally substituted

C_{1-6}-alkylthio, and halogen, where any C_{1-6}-alkyl as an amino substituent is optionally substituted with hydroxy, C_{1-6}-alkoxy, amino, mono- and di(C_{1-6}-alkyl)amino, carboxy, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylaminocarbonyl, or halogen(s); and

~~R^1, R^2, and R^3, when attached to a nitrogen atom, independently are selected from hydrogen, optionally substituted C_{1-6}-alkyl, hydroxy, optionally substituted C_{1-6}-alkoxy, optionally substituted C_{1-6}-alkoxycarbonyl, optionally substituted C_{1-6}-alkylcarbonyl, formyl, mono- and di(C_{1-6}-alkyl)-aminocarbonyl, amino, C_{1-6}-alkylcarbonylamino, mono- and di(C_{1-6}-alkyl)amino, C_{1-6}-alkylsulphonyl, and C_{1-6}-alkylsulphinyl; where any C_{1-6}-alkyl as an amino substituent is optionally substituted with hydroxy, C_{1-6}-alkoxy, amino, mono- and di(C_{1-6}-alkyl)amino, carboxy, C_{1-6}-alkyl-carbonylamino, C_{1-6}-alkylaminocarbonyl, or halogen(s), and wherein any aryl, heterocyclyl and heteroaryl may be optionally substituted;~~

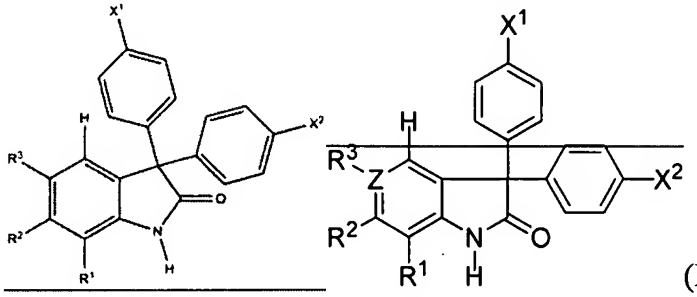
or wherein R^1 and R^2 together with the carbon ~~and/or nitrogen~~ atoms to which they are attached form a heterocyclic ring, a heteroaromatic ring, an aromatic ring or a carbocyclic ring;

~~Z is CH or N;~~ and

X^1 and X^2 are independently selected from halogen, OR^6, $OCOR^5$, $N(R^6)_2$, $NHCOR^5$, $NHSO_2R^5$, and $NHCON(R^6)_2$, wherein R^5 is selected from C_{1-6}-alkyl, optionally substituted aryl and optionally substituted heteroaryl, and each R^6 independently is selected from hydrogen, C_{1-6}-alkyl, optionally substituted aryl and optionally substituted heteroaryl; and

pharmaceutically acceptable salts and prodrugs thereof.

Claim 26 (**currently amended**): A method of treating a mammal suffering from or being susceptible to cancer, the method comprising administering to the mammal a therapeutically effective amount of a 3,3-diphenyl-1,3-dihydro-indol-2-one type compound of the formula (IIc)



(IIc)

wherein

R^1 is selected from hydrogen, halogen, C_{1-6}-alkyl, trifluoromethyl and C_{1-6}-alkoxy;

R^2 is selected from hydrogen, halogen, optionally substituted aryl, optionally substituted aryloxy, and optionally substituted heteroaryl;

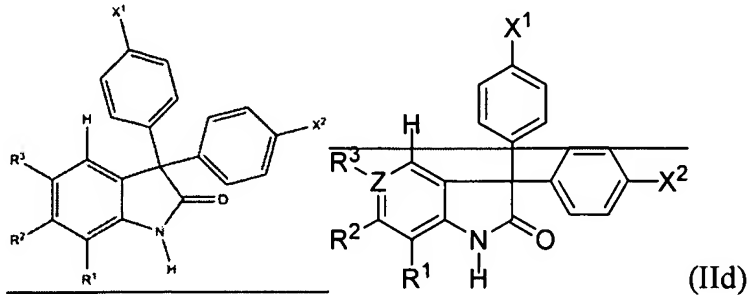
R^3 is selected from hydrogen, optionally substituted C_{1-6}-alkoxy, halogen, cyano, and optionally substituted aryl, optionally substituted aryloxy, optionally substituted heteroaryl, amino, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylsulphonylamino, and mono- and di(C_{1-6}-alkyl)aminosulfonyl;

~~Z is CH or N;~~ and

one of X^1 and X^2 is selected from halogen, OR^6, $OCOR^5$, $N(R^6)_2$, $NHCOR^5$, $NHSO_2R^5$, and $NHCON(R^6)_2$, wherein R^5 is selected from C_{1-6}-alkyl, optionally substituted aryl and optionally substituted heteroaryl, and each R^6 independently is selected from hydrogen, C_{1-6}-alkyl, optionally substituted aryl and optionally substituted heteroaryl; and the other of X^1 and X^2 is selected from optionally substituted C_{1-6}-alkyl, optionally substituted C_{2-6}-alkenyl, carboxy, optionally substituted C_{1-6}-alkoxycarbonyl, optionally substituted C_{1-6}-alkylcarbonyl, formyl, carbamoyl, mono- and di(C_{1-6}-alkyl)aminocarbonyl, cyano, aryl, arylcarbonyl, heterocyclyl, heterocyclylcarbonyl, heteroaryl, heteroarylcarbonyl, where any C_{1-6}-alkyl as an amino substituent is optionally substituted with hydroxy, C_{1-6}-alkoxy, amino, mono- and di(C_{1-6}-alkyl)amino, carboxy, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylaminocarbonyl, or halogen(s), and wherein any aryl, heterocyclyl and heteroaryl may be optionally substituted; and

pharmaceutically acceptable salts and prodrugs thereof.

Claim 27 (**currently amended**): A method of treating a mammal suffering from or being susceptible to cancer, the method comprising administering to the mammal a therapeutically effective amount of a 3,3-diphenyl-1,3-dihydro-indol-2-one type compound of the formula (IId)



(IId)

wherein

R^1, R^2, and R^3~~, when attached to a carbon atom,~~ independently are selected from hydrogen, optionally substituted C_{1-6}-alkyl, optionally substituted C_{2-6}-alkenyl, hydroxy, optionally substituted C_{1-6}-alkoxy, optionally substituted C_{2-6}-alkenyloxy, carboxy, optionally substituted C_{1-6}-alkoxycarbonyl, optionally substituted C_{1-6}-alkylcarbonyl, optionally substituted C_{1-6}-alkylcarbonyloxy, formyl, amino, mono- and di(C_{1-6}-alkyl)amino, carbamoyl, mono- and di(C_{1-6}-alkyl)aminocarbonyl, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylsulphonylamino, cyano, carbamido, mono- and di(C_{1-6}-alkyl)aminocarbonylamino, C_{1-6}-alkanoyloxy, C_{1-6}-alkylsulphonyl, C_{1-6}-alkylsulphinyl, aminosulfonyl, mono- and di(C_{1-6}-alkyl)aminosulfonyl, nitro, optionally substituted C_{1-6}-alkylthio, and halogen, where any C_{1-6}-alkyl as an amino substituent is optionally substituted with hydroxy, C_{1-6}-alkoxy, amino, mono- and di(C_{1-6}-alkyl)amino, carboxy, C_{1-6}-alkylcarbonylamino, C_{1-6}-alkylaminocarbonyl, or halogen(s); and

~~R^1, R^2, and R^3, when attached to a nitrogen atom, independently are selected from hydrogen, optionally substituted C_{1-6}-alkyl, hydroxy, optionally substituted C_{1-6}-alkoxy, optionally substituted C_{1-6}-alkoxycarbonyl, optionally substituted C_{1-6}-alkylcarbonyl, formyl, mono- and di(C_{1-6}-alkyl)-aminocarbonyl, amino, C_{1-6}-alkylcarbonylamino, mono- and di(C_{1-6}-alkyl)amino, C_{1-6}-alkylsulphonyl, and C_{1-6}-alkylsulphinyl; where any C_{1-6}-alkyl as an amino substituent is optionally substituted with hydroxy, C_{1-6}-alkoxy, amino, mono- and di(C_{1-6}-alkyl)amino, carboxy, C_{1-6}-alkyl-~~

~~carbonylamino, C_{1-6}-alkylaminocarbonyl, or halogen(s), and wherein any aryl, heterocyclyl and heteroaryl may be optionally substituted;~~

or wherein R^1 and R^2 together with the carbon ~~and/or nitrogen~~ atoms to which they are attached form a heterocyclic ring, a heteroaromatic ring, an aromatic ring or a carbocyclic ring;

~~Z is CH or N;~~ and

one of X^1 and X^2 is selected from halogen, OR^6, $OCOR^5$, $N(R^6)_2$, $NHCOR^5$, $NHSO_2R^5$, and $NHCON(R^6)_2$, wherein R^5 is selected from C_{1-6}-alkyl, optionally substituted aryl and optionally substituted heteroaryl, and each R^6 independently is selected from hydrogen, C_{1-6}-alkyl, optionally substituted aryl and optionally substituted heteroaryl; and the other of X^1 and X^2 is selected from optionally substituted C_{1-6}-alkyl, optionally substituted C_{2-6}-alkenyl, carboxy, optionally substituted C_{1-6}-alkoxycarbonyl, optionally substituted C_{1-6}-alkylcarbonyl, formyl, carbamoyl, mono- and di(C_{1-6}-alkyl)aminocarbonyl, cyano, aryl, arylcarbonyl, heterocyclyl, heterocyclylcarbonyl, heteroaryl, heteroarylcarbonyl, where any C_{1-6}-alkyl as an amino substituent is optionally substituted with hydroxy, C_{1-6}-alkoxy, amino, mono- and di(C_{1-6}-alkyl)amino, carboxy, C_{1-6}-alkyl-carbonylamino, C_{1-6}-alkylaminocarbonyl, or halogen(s), and wherein any aryl, heterocyclyl and heteroaryl may be optionally substituted; and

pharmaceutically acceptable salts and prodrugs thereof.

Claim 28 (**previously presented**): The method according to claim 1, wherein the compound is selected from Items 1 to 225 listed below:

1 5-Amino-6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one

2 5-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one

3 5-Fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

4 3,3-Bis-(4-hydroxy-phenyl)-5-nitro-1,3-dihydro-indol-2-one

5 3,3-Bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

6 6-Bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

7 6-Bromo-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

8 6-Bromo-3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one

9 6-Bromo-3,3-bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indole-5-carbonitrile

10 6-Bromo-3,3-bis-(4-hydroxy-phenyl)-5-methoxy-7-methyl-1,3-dihydro-indol-2-one

11 6-Bromo-3,3-bis-(4-hydroxy-phenyl)-7-methoxy-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one;

12 6-Bromo-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

13 6-Bromo-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one

14 6-Bromo-5-ethyl-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one

15 6-Bromo-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-5-carbonitrile

16 6-Bromo-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one

17 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

18 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

19 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one

20 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indole-5-carbonitrile

21 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-5-methoxy-7-methyl-1,3-dihydro-indol-2-one

22 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methoxy-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

23 6-Chloro-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

24 6-Chloro-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one

25 6-Chloro-5-ethyl-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one

26 6-Chloro-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-5-carbonitrile

27 6-Chloro-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one

28 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-5-methyl-7-methoxy-1,3-dihydro-indol-2-one;

29 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methoxy-2-oxo-2,3-dihydro-1H-indole-5-carbonitrile;

30 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methoxy-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one;

31 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methoxy-5-methyl-1,3-dihydro-indol-2-one;

32 6-Chloro-5-ethyl-3,3-bis-(4-hydroxy-phenyl)-7-methoxy-1,3-dihydro-indol-2-one;

33 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-5,7-dimethoxy-1,3-dihydro-indol-2-one;

34 N-{4-[3-(4-Acetylamino-phenyl)-5-chloro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-acetamide;

35 N-{4-[5-Chloro-3-(4-methanesulfonylamino-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-methanesulfonamide

36 N-{4-[3-(4-Acetylamino-phenyl)-6-chloro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-acetamide;

37 N-{4-[6-Chloro-3-(4-methanesulfonylamino-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-methanesulfonamide;

38 N-{4-[3-(4-Acetylamino-phenyl)-5-chloro-7-methoxy-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-acetamide;

39 N-{4-[5-Chloro-3-(4-methanesulfonylamino-phenyl)-7-methoxy-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-methanesulfonamide;

40 N-{4-[3-(4-Acetylamino-phenyl)-6-chloro-7-methoxy-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-acetamide; and

41 N-{4-[6-Chloro-3-(4-methanesulfonylamino-phenyl)-7-methoxy-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-methanesulfonamide

42 2-Chloro-6,6-bis-(4-hydroxy-phenyl)-3-methyl-4,6-dihydro-3H-pyrrolo[2,3-d]imidazol-5-one

43 Acetic acid 4-[6-(4-acetoxy-phenyl)-2-chloro-3-methyl-5-oxo-3,4,5,6-tetrahydro-pyrrolo[2,3-d]imidazol-6-yl]-phenyl ester

44 6,6-Bis-(4-amino-phenyl)-2-chloro-3-methyl-4,6-dihydro-3H-pyrrolo[2,3-d]imidazol-5-one

45 2-Chloro-6,6-bis-(4-dimethylamino-phenyl)-3-methyl-4,6-dihydro-3H-pyrrolo[2,3-d]imidazol-5-one

46 N-{4-[6-(4-Acetylamino-phenyl)-2-chloro-3-methyl-5-oxo-3,4,5,6-tetrahydro-pyrrolo[2,3-d]imidazol-6-yl]-phenyl}-acetamide

47 N-{4-[2-Chloro-6-(4-methanesulfonylamino-phenyl)-3-methyl-5-oxo-3,4,5,6-tetrahydro-pyrrolo[2,3-d]imidazol-6-yl]-phenyl}-methanesulfonamide

48 4,4-Bis-(4-hydroxy-phenyl)-1-methyl-4,6-dihydro-1H-pyrrolo[2,3-c]pyrazol-5-one

49 Acetic acid 4-[4-(4-acetoxy-phenyl)-1-methyl-5-oxo-1,4,5,6-tetrahydro-pyrrolo[2,3-c]pyrazol-4-yl]-phenyl ester

50 4,4-Bis-(4-amino-phenyl)-1-methyl-4,6-dihydro-1H-pyrrolo[2,3-c]pyrazol-5-one

51 N-{4-[4-(4-Methanesulfonylamino-phenyl)-1-methyl-5-oxo-1,4,5,6-tetrahydro-pyrrolo[2,3-c]pyrazol-4-yl]-phenyl}-methanesulfonamide

52 4,4-Bis-(4-dimethylamino-phenyl)-1-methyl-4,6-dihydro-1H-pyrrolo[2,3-c]pyrazol-5-one

53 N-{4-[4-(4-Acetylamino-phenyl)-1-methyl-5-oxo-1,4,5,6-tetrahydro-pyrrolo[2,3-c]pyrazol-4-yl]-phenyl}-acetamide

54 4,4-Bis-(4-hydroxy-phenyl)-2-methyl-2,6-dihydro-4H-pyrrolo[2,3-c]pyrazol-5-one

55 Acetic acid 4-[4-(4-acetoxy-phenyl)-2-methyl-5-oxo-2,4,5,6-tetrahydro-pyrrolo[2,3-c]pyrazol-4-yl]-phenyl ester

56 4,4-Bis-(4-amino-phenyl)-2-methyl-2,6-dihydro-4H-pyrrolo[2,3-c]pyrazol-5-one

57 4,4-Bis-(4-dimethylamino-phenyl)-2-methyl-2,6-dihydro-4H-pyrrolo[2,3-c]pyrazol-5-one

58 N-{4-[4-(4-Acetylamino-phenyl)-2-methyl-5-oxo-2,4,5,6-tetrahydro-pyrrolo[2,3-c]pyrazol-4-yl]-phenyl}-acetamide

59 N-{4-[4-(4-Methanesulfonylamino-phenyl)-2-methyl-5-oxo-2,4,5,6-tetrahydro-pyrrolo[2,3-c]pyrazol-4-yl]-phenyl}-methanesulfonamide

60 4,4-Bis-(4-hydroxy-phenyl)-4,6-dihydro-thieno[2,3-b]pyrrol-5-one

61 Acetic acid 4-[4-(4-acetoxy-phenyl)-5-oxo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yl]-phenyl ester

62 4,4-Bis-(4-amino-phenyl)-4,6-dihydro-thieno[2,3-b]pyrrol-5-one

63 4,4-Bis-(4-dimethylamino-phenyl)-4,6-dihydro-thieno[2,3-b]pyrrol-5-one

64 N-{4-[4-(4-Acetylamino-phenyl)-5-oxo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yl]-phenyl}-acetamide

65 N-{4-[4-(4-Methanesulfonylamino-phenyl)-5-oxo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yl]-phenyl}-methanesulfonamide

66 2-Chloro-4,4-bis-(4-hydroxy-phenyl)-4,6-dihydro-thieno[2,3-b]pyrrol-5-one

67 Acetic acid 4-[4-(4-acetoxy-phenyl)-2-chloro-5-oxo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yl]-phenyl ester

68 4,4-Bis-(4-amino-phenyl)-2-chloro-4,6-dihydro-thieno[2,3-b]pyrrol-5-one

69 2-Chloro-4,4-bis-(4-dimethylamino-phenyl)-4,6-dihydro-thieno[2,3-b]pyrrol-5-one

70 N-{4-[4-(4-Acetylamino-phenyl)-2-chloro-5-oxo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yl]-phenyl}-acetamide

71 N-{4-[2-Chloro-4-(4-methanesulfonylamino-phenyl)-5-oxo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yl]-phenyl}-methanesulfonamide

72 4,4-Bis-(4-hydroxy-phenyl)-4,6-dihydro-furo[2,3-b]pyrrol-5-one

73 Acetic acid 4-[4-(4-acetoxy-phenyl)-5-oxo-5,6-dihydro-4H-furo[2,3-b]pyrrol-4-yl]-phenyl ester

74 4,4-Bis-(4-amino-phenyl)-4,6-dihydro-furo[2,3-b]pyrrol-5-one

75 4,4-Bis-(4-dimethylamino-phenyl)-4,6-dihydro-furo[2,3-b]pyrrol-5-one

76 N-{4-[4-(4-Acetylamino-phenyl)-5-oxo-5,6-dihydro-4H-furo[2,3-b]pyrrol-4-yl]-phenyl}-acetamide

77 N-{4-[4-(4-Methanesulfonylamino-phenyl)-5-oxo-5,6-dihydro-4H-furo[2,3-b]pyrrol-4-yl]-phenyl}-methanesulfonamide

78 2-Chloro-4,4-bis-(4-hydroxy-phenyl)-4,6-dihydro-furo[2,3-b]pyrrol-5-one

79 Acetic acid 4-[4-(4-acetoxy-phenyl)-2-chloro-5-oxo-5,6-dihydro-4H-furo[2,3-b]pyrrol-4-yl]-phenyl ester

80 4,4-Bis-(4-amino-phenyl)-2-chloro-4,6-dihydro-furo[2,3-b]pyrrol-5-one

81 2-Chloro-4,4-bis-(4-dimethylamino-phenyl)-4,6-dihydro-furo[2,3-b]pyrrol-5-one

82 N-{4-[4-(4-Acetylamino-phenyl)-2-chloro-5-oxo-5,6-dihydro-4H-furo[2,3-b]pyrrol-4-yl]-phenyl}-acetamide

83 N-{4-[2-Chloro-4-(4-methanesulfonylamino-phenyl)-5-oxo-5,6-dihydro-4H-furo[2,3-b]pyrrol-4-yl]-phenyl}-methanesulfonamide

84 3,3-Bis-(4-hydroxy-phenyl)-6-methyl-3,8-dihydro-1H-1,8-diaza-as-indacen-2-one

85 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-methyl-2-oxo-1,2,3,8-tetrahydro-1,8-diaza-as-indacen-3-yl]-phenyl ester

86 3,3-Bis-(4-amino-phenyl)-6-methyl-3,8-dihydro-1H-1,8-diaza-as-indacen-2-one

87 3,3-Bis-(4-dimethylamino-phenyl)-6-methyl-3,8-dihydro-1H-1,8-diaza-as-indacen-2-one

88 N-{4-[3-(4-Acetylamino-phenyl)-6-methyl-2-oxo-1,2,3,8-tetrahydro-1,8-diaza-as-indacen-3-yl]-phenyl}-acetamide

89 N-{4-[3-(4-Methanesulfonylamino-phenyl)-6-methyl-2-oxo-1,2,3,8-tetrahydro-1,8-diaza-as-indacen-3-yl]-phenyl}-methanesulfonamide

90 3,3-Bis-(4-hydroxy-phenyl)-1,3-dihydro-benzo[g]indol-2-one

91 Acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-benzo[g]indol-3-yl]-phenyl ester

92 3,3-Bis-(4-amino-phenyl)-1,3-dihydro-benzo[g]indol-2-one

93 3,3-Bis-(4-dimethylamino-phenyl)-1,3-dihydro-benzo[g]indol-2-one

94 N-{4-[3-(4-Acetylamino-phenyl)-2-oxo-2,3-dihydro-1H-benzo[g]indol-3-yl]-phenyl}-acetamide

95 N-{4-[3-(4-Methanesulfonylamino-phenyl)-2-oxo-2,3-dihydro-1H-benzo[g]indol-3-yl]-phenyl}-methanesulfonamide

96 1-Amino-6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one

97 Acetic acid 4-[3-(4-acetoxy-phenyl)-1-amino-6-chloro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester

98 N-{4-[3-(4-Acetylamino-phenyl)-1-amino-6-chloro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-acetamide

99 N-{4-[1-Amino-6-chloro-3-(4-methanesulfonylamino-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-methanesulfonamide

100 Acetic acid 4-[3-(4-acetoxy-phenyl)-1-acetylamino-6-chloro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester

101 N-[3,3-Bis-(4-amino-phenyl)-6-chloro-7-methyl-2-oxo-2,3-dihydro-indol-1-yl]-acetamide

102 N-[6-Chloro-3,3-bis-(4-dimethylamino-phenyl)-7-methyl-2-oxo-2,3-dihydro-indol-1-yl]-acetamide

103 N-[3,3-Bis-(4-acetylamino-phenyl)-6-chloro-7-methyl-2-oxo-2,3-dihydro-indol-1-yl]-acetamide

104 N-[6-Chloro-3,3-bis-(4-methanesulfonylamino-phenyl)-7-methyl-2-oxo-2,3-dihydro-indol-1-yl]-acetamide

105 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indole-2-thione

106 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-7-methyl-2-thioxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester

107 3,3-Bis-(4-amino-phenyl)-6-chloro-7-methyl-1,3-dihydro-indole-2-thione

108 6-Chloro-3,3-bis-(4-dimethylamino-phenyl)-7-methyl-1,3-dihydro-indole-2-thione

109 N-{4-[3-(4-Acetylamino-phenyl)-6-chloro-7-methyl-2-thioxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-acetamide

110 Methanesulfonic acid 4-[6-chloro-3-(4-methanesulfonyloxy-phenyl)-7-methyl-2-thioxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester

111 Acetic acid 4-[4-(4-acetoxy-phenyl)-2-chloro-5-thioxo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yl]-phenyl ester

112 Acetic acid 4-[4-(4-acetoxy-phenyl)-2-chloro-5-thioxo-5,6-dihydro-4H-furo[2,3-b]pyrrol-4-yl]-phenyl ester

113 6,6-Bis-(4-amino-phenyl)-2-chloro-3-methyl-4,6-dihydro-thieno[3,2-b]pyrrole-5-thione

114 2-Chloro-6,6-bis-(4-dimethylamino-phenyl)-3-methyl-4,6-dihydro-3H-pyrrolo[2,3-d]imidazole-5-thione

115 N-{4-[6-(4-Acetylamino-phenyl)-3-chloro-5-thioxo-1,4,5,6-tetrahydro-pyrrolo[3,2-c]pyrazol-6-yl]-phenyl}-acetamide

116 Methanesulfonic acid 4-[2-chloro-4-(4-methanesulfonyloxy-phenyl)-5-thioxo-5,6-dihydro-4H-furo[2,3-b]pyrrol-4-yl]-phenyl ester

117 6-Chloro-7-cyclopropyl-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

118 6-Chloro-7-cyclopropyl-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

119 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-trifluoromethyl-1,3-dihydro-indol-2-one

120 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-trifluoromethyl-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

121 6-Chloro-7-cyclopropoxy-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

122 6-Chloro-7-cyclopropoxy-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

123 6-(4-Fluoro-phenoxy)-3,3-bis-(4-hydroxy-phenyl)-7-trifluoromethyl-1,3-dihydro-indol-2-one

124 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-7-cyclopropyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester

125 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-7-cyclopropyl-2-oxo-2,3-dihydro-1H-pyrrolo[3,2-c]pyridin-3-yl]-phenyl ester

126 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-2-oxo-7-trifluoromethyl-2,3-dihydro-1H-indol-3-yl]-phenyl ester

127 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-2-oxo-7-trifluoromethyl-2,3-dihydro-1H-pyrrolo[3,2-c]pyridin-3-yl]-phenyl ester

128 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-7-cyclopropoxy-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester

129 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-7-cyclopropoxy-2-oxo-2,3-dihydro-1H-pyrrolo[3,2-c]pyridin-3-yl]-phenyl ester

130 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-(4-fluoro-phenoxy)-2-oxo-7-trifluoromethyl-2,3-dihydro-1H-indol-3-yl]-phenyl ester

131 Dimethylamino-acetic acid 4-{6-chloro-7-cyclopropyl-3-[4-(2-dimethylamino-acetoxy)-phenyl]-2-oxo-2,3-dihydro-1H-indol-3-yl}-phenyl ester

132 Dimethylamino-acetic acid 4-{6-chloro-7-cyclopropyl-3-[4-(2-dimethylamino-acetoxy)-phenyl]-2-oxo-2,3-dihydro-1H-pyrrolo[3,2-c]pyridin-3-yl}-phenyl ester

133 Dimethylamino-acetic acid 4-{6-chloro-3-[4-(2-dimethylamino-acetoxy)-phenyl]-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl}-phenyl ester

134 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-trifluoromethoxy-1,3-dihydro-indol-2-one

135 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-2-oxo-7-trifluoromethoxy-2,3-dihydro-1H-indol-3-yl]-phenyl ester

136 Dimethylamino-acetic acid 4-{6-chloro-3-[4-(2-dimethylamino-acetoxy)-phenyl]-2-oxo-7-trifluoromethoxy-2,3-dihydro-1H-indol-3-yl}-phenyl ester

137 6-Chloro-4-fluoro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one

138 3-Chloro-7,7-bis-(4-hydroxy-phenyl)-4-methyl-5,7-dihydro-pyrrolo[3,2-c]pyridazin-6-one

139 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-4-fluoro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester

140 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-4,7-dimethyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester

141 Acetic acid 4-[7-(4-acetoxy-phenyl)-3-chloro-4-methyl-6-oxo-6,7-dihydro-5H-pyrrolo[3,2-c]pyridazin-7-yl]-phenyl ester

142 6-Chloro-4,5-difluoro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one

143 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-4,5-difluoro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester

144 3,3-Bis-(4-hydroxy-phenyl)-3,6,7,8-tetrahydro-1H-1-aza-as-indacen-2-one

145 3,3-Bis-(4-hydroxy-phenyl)-1,3,6,7,8,9-hexahydro-benzo[g]indol-2-one

146 3,3-Bis-(4-hydroxy-phenyl)-7-trifluoromethyl-1,3-dihydro-indol-2-one

147 7-Chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

148 3,3-Bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-7-carbonitrile

149 7-Ethyl-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

150 3,3-Bis-(4-hydroxy-phenyl)-7-morpholin-4-yl-1,3-dihydro-indol-2-one

151 3,3-Bis-(4-hydroxy-phenyl)-7-isopropyl-1,3-dihydro-indol-2-one

152 7-tert-Butyl-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

153 3,3-Bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-7-carboxylic acid dimethylamide

154 3,3-Bis-(4-hydroxy-phenyl)-7-(4-methyl-piperazine-1-carbonyl)-1,3-dihydro-indol-2-one

155 3,3-Bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-5-carboxylic acid

156 3,3-Bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-5-carboxylic acid dimethylamide

157 3,3-Bis-(4-hydroxy-phenyl)-5-(morpholine-4-carbonyl)-1,3-dihydro-indol-2-one

158 3,3-Bis-(4-hydroxy-phenyl)-4-methoxy-1,3-dihydro-indol-2-one

159 3,3-Bis-(4-hydroxy-phenyl)-6-methoxy-1,3-dihydro-indol-2-one

160 3,3-Bis-(4-hydroxy-phenyl)-5-(4-methyl-piperazine-1-carbonyl)-1,3-dihydro-indol-2-one

161 6-Chloro-3,3-bis-(4-mercapto-phenyl)-7-methyl-1,3-dihydro-indol-2-one

162 N-{4-[3-(4-Acetylamino-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-acetamide

163 3,3-Bis-(4-hydroxy-phenyl)-7-(3-methoxy-prop-1-ynyl)-1,3-dihydro-indol-2-one

164 3,3-Bis-(4-hydroxy-phenyl)-7-pyridin-3-yl-1,3-dihydro-indol-2-one

165 7-Bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

166 6-Chloro-3,3-bis-(4-methanesulfonyl-phenyl)-7-methyl-1,3-dihydro-indol-2-one

167 6,6-Bis-(4-hydroxy-phenyl)-4,6-dihydro-pyrrolo[3,2-d]thiazol-5-one

168 6,6-Bis-(4-hydroxy-phenyl)-2-methyl-4,6-dihydro-pyrrolo[3,2-d]thiazol-5-one

169 6,6-Bis-(4-hydroxy-phenyl)-2-isopropyl-4,6-dihydro-pyrrolo[3,2-d]thiazol-5-one

170 2-Chloro-6,6-bis-(4-hydroxy-phenyl)-4,6-dihydro-pyrrolo[3,2-d]thiazol-5-one

171 4,4-Bis-(4-hydroxy-phenyl)-4,6-dihydro-pyrrolo[3,2-d]isothiazol-5-one

172 3,3-Bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-pyrrolo[2,3-c]pyridin-2-one

173 3,3-Bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-pyrrolo[3,2-b]pyridin-2-one

174 3,3-Bis-(4-fluoro-phenyl)-7-methyl-1,3-dihydro-pyrrolo[3,2-b]pyridin-2-one

175 3,3-Bis-(4-fluoro-phenyl)-7-methyl-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

176 3,3-Bis-(4-fluoro-phenyl)-7-isopropyl-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one

177 3,3-Bis-(4-hydroxy-phenyl)-3,6,7,8-tetrahydro-1H-1,5-diaza-as-indacen-2-one

178 3,3-Bis-(4-hydroxy-phenyl)-3,6,7,8-tetrahydro-1H-1,4-diaza-as-indacen-2-one

179 3,3-Bis-(4-hydroxy-phenyl)-1,3,6,7,8,9-hexahydro-pyrrolo[3,2-c]quinolin-2-one

180 3,3-Bis-(4-hydroxy-phenyl)-1,3,6,7,8,9-hexahydro-pyrrolo[3,2-c]isoquinolin-2-one

181 5-Fluoro-3,3-bis-(4-hydroxy-phenyl)-3,6,7,8-tetrahydro-1H-1-aza-as-indacen-2-one

182 7-Ethyl-5-fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

183 3,3-Bis-(4-hydroxy-phenyl)-1,3,6,8-tetrahydro-7-oxa-1-aza-as-indacen-2-one

184 3,3-Bis-(4-hydroxy-phenyl)-1,3,7,8-tetrahydro-6-oxa-1-aza-as-indacen-2-one

185 3,3-Bis-(4-hydroxy-phenyl)-1,6,7,9-tetrahydro-3H-8-oxa-1-aza-cyclopenta[a]naphthalen-2-one

186 3,3-Bis-(4-hydroxy-phenyl)-1,7,8,9-tetrahydro-3H-pyrano[2,3-g]indol-2-one

187 3,3-Bis-(4-hydroxy-phenyl)-7-methyl-3,6,7,8-tetrahydro-1H-1,7-diaza-as-indacen-2-one

188 3,3-Bis-(4-hydroxy-phenyl)-7-methyl-1,3,7,8-tetrahydro-1,7-diaza-as-indacene-2,6-dione

189 3,3-Bis-(4-hydroxy-phenyl)-7,8,8-trimethyl-1,3,7,8-tetrahydro-1,7-diaza-as-indacene-2,6-dione

190 3,3-Bis-(4-hydroxy-phenyl)-5-iodo-1,3-dihydro-indol-2-one

191 5-Amino-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

192 5-Amino-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one

193 6-Bromo-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one

194 7-Fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

195 3,3-Bis-(4-hydroxy-phenyl)-7-methoxy-1,3-dihydro-indol-2-one

196 4,7-Dichloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

197 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-1,7-dimethyl-1,3-dihydro-indol-2-one

198 6-Chloro-3,3-bis-(4-fluoro-phenyl)-7-methyl-1,3-dihydro-indol-2-one

199 3,3-Bis-(4-hydroxy-phenyl)-7-(morpholine-4-carbonyl)-1,3-dihydro-indol-2-one

200 3,3-Bis-(4-hydroxy-phenyl)-1,3-dihydro-pyrrolo[2,3-d]pyridin-2-one

201 N-{4-[6-Chloro-3-(4-methanesulfonylamino-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-methanesulfonamide

202 3,3-Bis-(4-hydroxy-phenyl)-4,7-dimethyl-1,3-dihydro-indol-2-one

203 3,3-Bis-(4-hydroxy-phenyl)-7-iodo-1,3-dihydro-indol-2-one

204 3,3-Bis-(4-hydroxy-phenyl)-7-pyridin-4-yl-1,3-dihydro-indol-2-one

205 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester

206 3,3-Bis-(4-hydroxy-phenyl)-5-phenyl-1,3-dihydro-indol-2-one

207 3,3-Bis-(4-hydroxy-phenyl)-7-thiophen-2-yl-1,3-dihydro-indol-2-one

208 3,3-Bis-(4-hydroxy-phenyl)-5-pyridin-4-yl-1,3-dihydro-indol-2-one

209 3,3-Bis-(4-hydroxy-phenyl)-5-thiophen-2-yl-1,3-dihydro-indol-2-one

210 5,7-Difluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

211 6-Fluoro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one

212 3,3-Bis-(4-hydroxy-phenyl)-6-methoxy-7-methyl-1,3-dihydro-indol-2-one

213 6,7-Difluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

214 6-Chloro-7-fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

215 5-Fluoro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one

216 3,3-Bis-(4-hydroxy-phenyl)-5-methoxy-7-methyl-1,3-dihydro-indol-2-one

217 3,3-Bis-(4-hydroxy-phenyl)-1,3-dihydro-pyrrolo[2,3-b]pyridin-2-one

218 7-Chloro-3,3-bis-(4-hydroxy-phenyl)-4-methoxy-1,3-dihydro-indol-2-one

219 6-Fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one

220 N-[3,3-Bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-indol-1-yl]-acetamide

221 5-[3,3-Bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-6-yloxy]-pentanoic acid methyl ester

222 5-[3,3-Bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-6-yloxy]-pentanoic acid

223 5-[3,3-Bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-5-yloxy]-pentanoic acid methyl ester

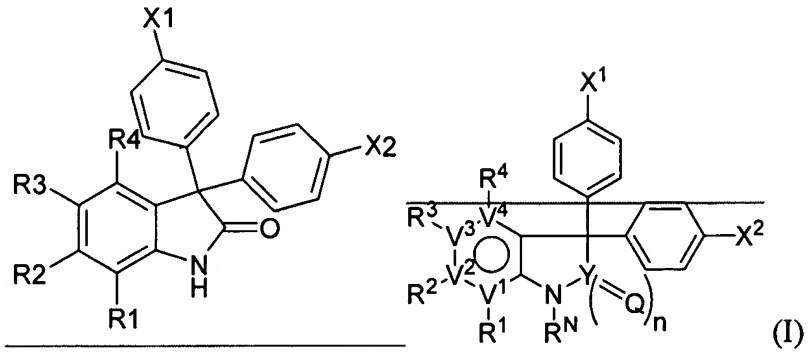
224 5-[3,3-Bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-5-yloxy]-pentanoic acid

225 7-Chloro-6-fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one.

Claim 29 (**previously presented**): The method according to claim 1, wherein the medicament further comprises one or more other chemotherapeutic agents.

Claim 30 (**canceled**).

Claim 31 (**currently amended**): A compound of the general formula (I)



(I)

as defined in claim 1, with the proviso that the compound is not one selected from

3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one,

3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-4,5-dimethyl-1,3-dihydro-indol-2-one ;

3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;

5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;

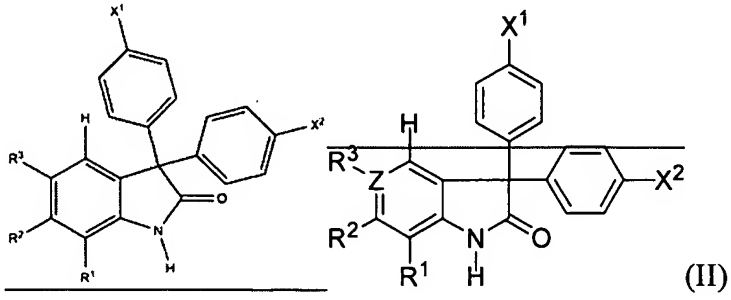
3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester; and

acetic acid 4-[3-(4-acetoxy-phenyl)-5-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester.

Claim 32 (**currently amended**): A 3,3-Diphenyl-1,3-dihydro-indol-2-one type compound of the formula (II)



(II)

as defined in claim 24, with the proviso that the compound is not one selected from:

3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one,

3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-4,5-dimethyl-1,3-dihydro-indol-2-one ;

3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;

5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

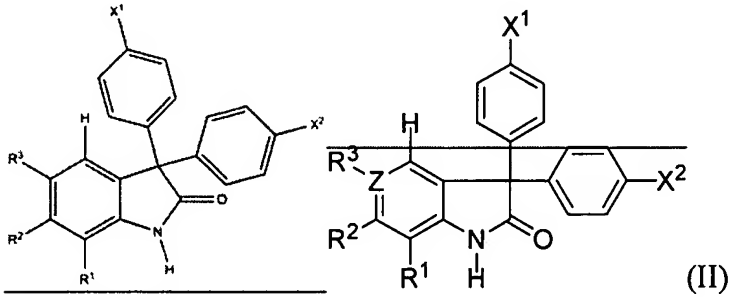
6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester; and

acetic acid 4-[3-(4-acetoxy-phenyl)-5-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester.

Claim 33 (**previously presented**): A pharmaceutical composition comprising a compound as defined in claim 1 and a pharmaceutically acceptable carrier.

Claim 34 (**currently amended**): A 3,3-Diphenyl-1,3-dihydro-indol-2-one type compound of the formula (II)



(II)

as defined in claim 25, with the proviso that the compound is not one selected from:

3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one,

3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-4,5-dimethyl-1,3-dihydro-indol-2-one ;

3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;

5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;

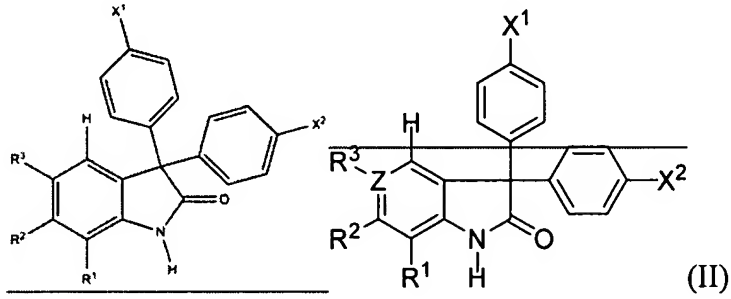
3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester; and

acetic acid 4-[3-(4-acetoxy-phenyl)-5-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester.

Claim 35 (**currently amended**): A 3,3-Diphenyl-1,3-dihydro-indol-2-one type compound of the formula (II)



(II)

as defined in claim 26, with the proviso that the compound is not one selected from:

3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one,

3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-4,5-dimethyl-1,3-dihydro-indol-2-one ;

3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;

5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;

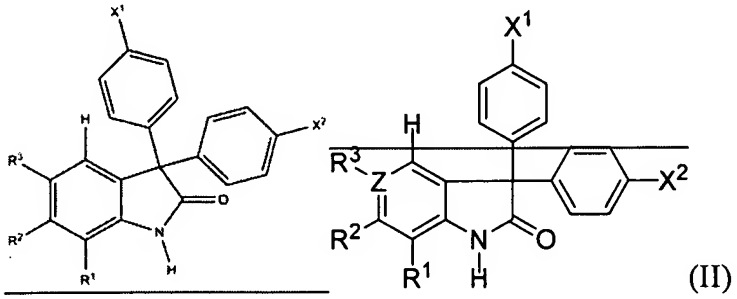
3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester; and

acetic acid 4-[3-(4-acetoxy-phenyl)-5-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester.

Claim 36 (**currently amended**): A 3,3-Diphenyl-1,3-dihydro-indol-2-one type compound of the formula (II)



(II)

as defined in claim 27, with the proviso that the compound is not one selected from:

3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one,

3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-4,5-dimethyl-1,3-dihydro-indol-2-one ;

3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;

5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;

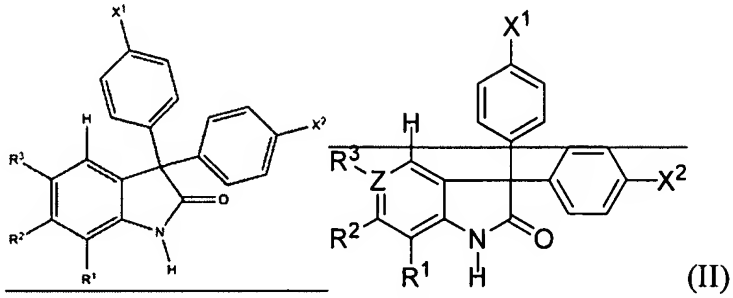
3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester; and

acetic acid 4-[3-(4-acetoxy-phenyl)-5-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester.

Claim 37 (**currently amended**): A 3,3-Diphenyl-1,3-dihydro-indol-2-one type compound of the formula (II)



(II)

as defined in claim 28, with the proviso that the compound is not one selected from:

3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one,

3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-4,5-dimethyl-1,3-dihydro-indol-2-one ;

3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;

5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester; and

acetic acid 4-[3-(4-acetoxy-phenyl)-5-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester.

Claim 38 (**new**): The method according to claim 1, wherein both of X^1 and X^2 are hydroxyl.